

**REMARKS**

Further consideration of this application is requested. Claims 8-11 remain active in the application. Of these, claim 10 has been allowed while claims 8, 9 and 10 are rejected over U.S. patent 3,919,316 to Molloy. Claim 8 is amended to exclude the compound of Molloy by adding a further "proviso" to claim 8, namely item (e).

Applicants submit that the claims of this application are now in condition for allowance and that claims 8, 9 and 11 should also be allowed with claim 10.

Favorable action is solicited.

Attached hereto is a marked-up version of the changes made to the claims by the current amendment. The attached page is captioned "**Version With Markings To Show Changes Made.**"

Respectfully submitted,

**NIXON & VANDERHYE P.C.**

By: \_\_\_\_\_



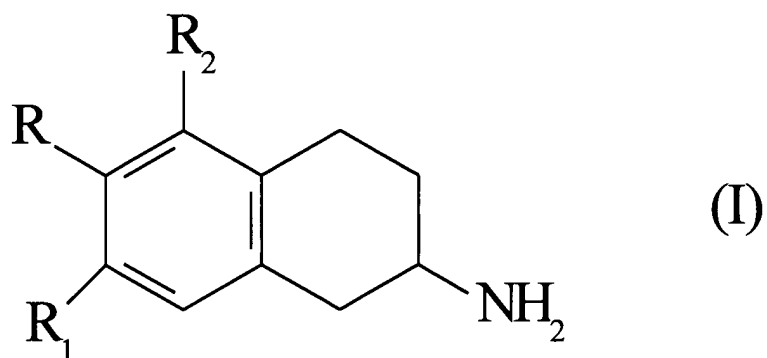
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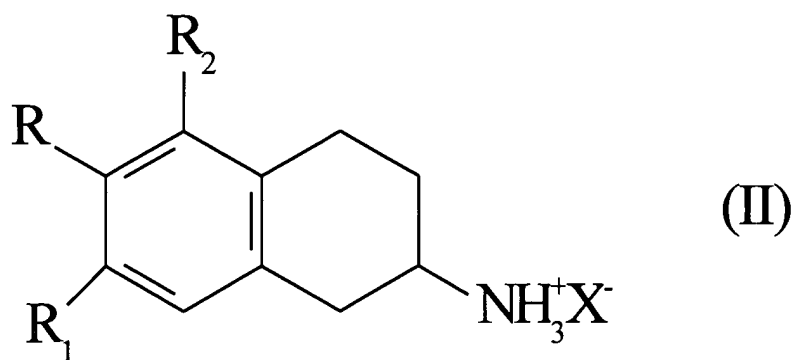
VERSION WITH MARKINGS TO SHOW CHANGES MADE

IN THE CLAIMS

8. (Twice Amended) A 2-aminoteraline of the formula (I)



or a pharmacologically acceptable salt of the formula (II)



wherein:

R and R<sub>1</sub> are independently halogen, hydroxy, or C<sub>1</sub>-C<sub>4</sub> alkoxy optionally substituted in position ω with a group selected from OH, NH<sub>2</sub> or NR<sub>3</sub>R<sub>4</sub>, wherein R<sub>3</sub> and R<sub>4</sub> are independently H, C<sub>1</sub>-C<sub>4</sub> alkyl, unsubstituted or substituted in position ω with groups OH, NH<sub>2</sub>, C<sub>1</sub>-C<sub>4</sub> alkanoyl, C<sub>1</sub>-C<sub>4</sub> alkyl, carbamoyl, carbamoyloxy, amino, or amino-substituted NR<sub>3</sub>R<sub>4</sub>, where R<sub>3</sub> and R<sub>4</sub> have the above meanings,

R<sub>2</sub> is hydrogen, halogen, hydroxy or methoxy,  
with the proviso that the 2-aminotetraline excludes (a) R=R<sub>1</sub>=CH<sub>3</sub>O or OH, R<sub>2</sub>=H, (b) R=F, R<sub>1</sub>=CH<sub>3</sub>O or OH, R<sub>2</sub>=H, (c) R<sub>1</sub>=-OCH<sub>3</sub>, R=CH<sub>3</sub> and R<sub>2</sub>=H [or] (d) R=R<sub>1</sub>=R<sub>2</sub>=CH<sub>3</sub>O or (e) R=R<sub>1</sub>=Cl and R<sub>2</sub>=H,

and X<sup>-</sup> is the monovalent anion of a pharmacologically acceptable acid.